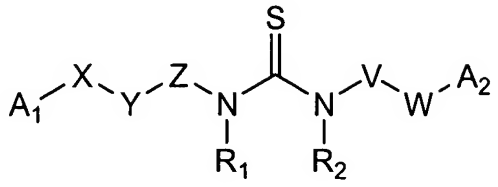
WHAT IS CLAIMED IS:

1. A compound of Formula 1



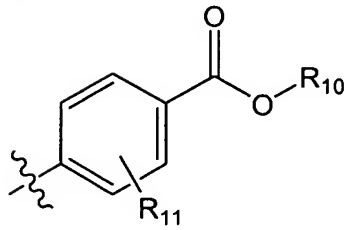
Formula 1

or a pharmaceutically acceptable salt thereof, wherein

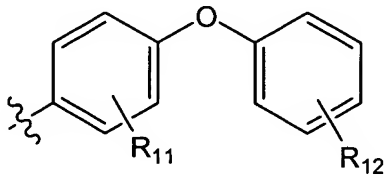
A_1 is an optionally substituted di-alkylamino, an optionally substituted aryl group, an optionally substituted 5- or 6- membered heteroaryl group, an optionally substituted bicyclic heteroaryl group having a 5-membered heteroaryl ring fused to a phenyl ring, an optionally substituted partially unsaturated or aromatic heterocyclic group having two 6-membered rings, an optionally substituted 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, an optionally substituted partially unsaturated 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a 5- or 6-membered heterocycloalkyl group fused to a phenyl or heteroaryl ring, or a fused or spiro 8 to 11-membered bicyclic heterocycloalkyl group containing at least one nitrogen atom and 0 to 3 additional heteroatoms;

A_2 is

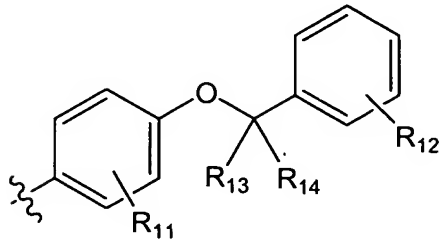
(i)



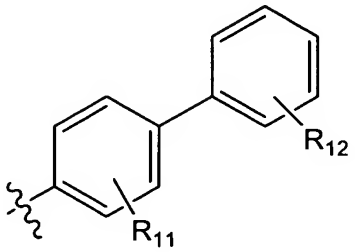
(ii)



(iii)



(iv)



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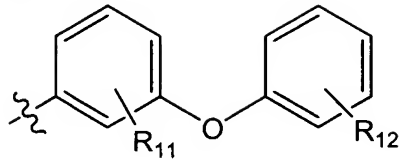
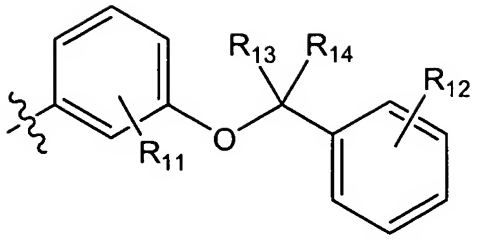


(vi)



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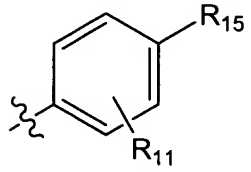
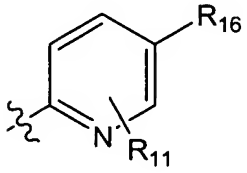
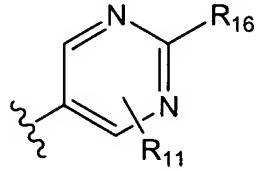


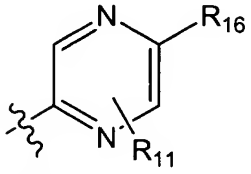
(viii)



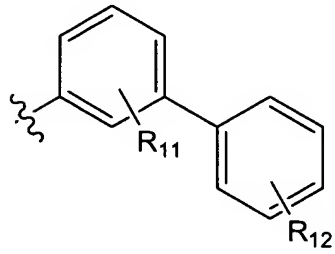
(ix)



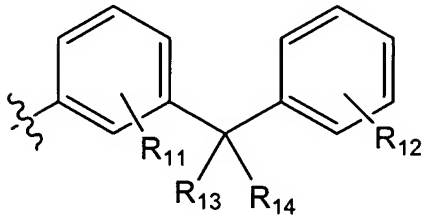
(x)



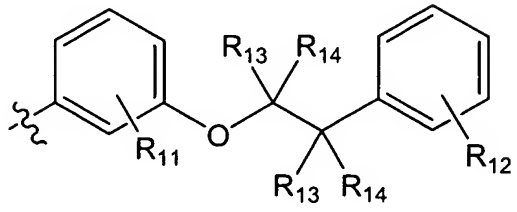
(xi)



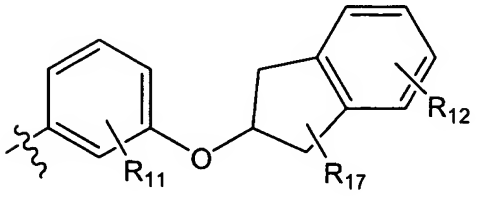
(xii)



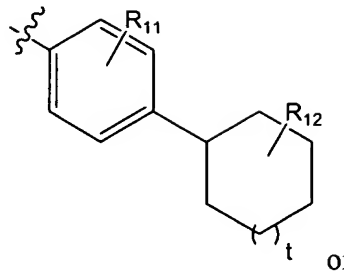
(xiii)



(xiv)

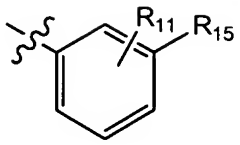


(xv)



or

(xvi)



;

t is 0 or 1;

X and W are independently O, S, NR, or absent, where R is hydrogen, optionally substituted C_1-C_6alkyl, or optionally substituted aryl(C_0-C_4alkyl);

V is C_1-C_6 alkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent;

Y is C_1-C_6 alkyl, C_1-C_6 alkyl substituted with C_3-C_7cycloalkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent;

wherein when V is absent, W is absent;

Z is carbonyl, thiocarbonyl, imino, or C_1-C_6alkylimino;

R_1 and R_2 are independently hydrogen, or

R_1 and R_2 are independently C_1-C_6alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, or

R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy;

R_{10} is C_1-C_6alkyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_6alkyl, C_1-C_6alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_2-C_6alkanoyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl; and

R_{13} and R_{14} are independently chosen at each occurrence from hydrogen and C_1-C_4alkyl;

R_{15} is C_4-C_6alkoxy or C_4-C_6alkyl;

R_{16} is C_2-C_6alkoxy or C_2-C_6alkyl; and

R_{17} represents 0 to 2 substituents independently chosen from halogen, methyl, and methoxy;

2. A compound or salt according to Claim 1 wherein

A_1 is a di-(C_1-C_6alkyl)amino, an aryl group, a 5- or 6- membered heteroaryl group, a bicyclic heteroaryl group having a 5-membered heteroaryl ring fused to a phenyl ring, a partially unsaturated or aromatic heterocyclic group having two 6-membered rings, a 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a partially unsaturated 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a 5- or 6-membered heterocycloalkyl group fused to a phenyl or heteroaryl ring, or a fused or spiro 8 to 11-membered bicyclic heterocycloalkyl group containing at least one nitrogen atom and 0 to 3 additional heteroatoms; each of which A_1 is substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy,

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) –GR_a where

G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl;

X and W are independently O, S, NR, or absent,

where R is hydrogen or R is C_1-C_6alkyl or aryl(C_0-C_4alkyl), each of which is substituted with 0 to 5 substitutents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, C_1-C_6alkyl, C_1-C_6alkoxy, and mono- and di-(C_1-C_6alkyl)amino;

V is independently C_1-C_6 alkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent;

Y is C_1-C_6 alkyl, C_1-C_6 alkyl substituted with C_3-C_7cycloalkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent;

wherein when V is absent, W is absent;

Z is carbonyl, thiocarbonyl, or imino; and

R_1 and R_2 are independently hydrogen, or

R_1 and R_2 are independently C_1-C_6alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, or

R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

3. A compound or salt according to Claim 1 or 2 in which Z is thiocarbonyl.

4. A compound or salt according to Claim 1 or 2 in which Z is imino or C_1-C_6alkylimino.

5. A compound or salt according to Claim 4 in which Z is imino or methylimino.

6. A compound or salt according to Claim 1 or 2 in which Z is carbonyl.

7. A compound or salt according to any one of Claims 1 to 6 in which X is oxygen and Y is -CH_2-.

8. A compound or salt according to any one of Claims 1 to 6 in which X is oxygen and Y is $-CH_2CH_2-$.

9. A compound or salt according to any one of Claims 1 to 6 wherein when X and Y are absent.

10. A compound or salt according to any one of Claims 1 to 9 wherein V and W are absent.

11. A compound or salt according to any one of Claims 1 to 9 in which V is C_1-C_2alkyl and W is absent.

12. A compound or salt according to any one of Claims 1 to 11 in which R_1 and R_2 are independently

hydrogen, or

C_1-C_4alkyl, C_2-C_4alkenyl, or C_2-C_4alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

13. A compound or salt according to Claim 12 in which R_1 and R_2 are independently hydrogen, methyl, or ethyl.

14. A compound or salt according to Claim 13 in which R_1 and R_2 are both hydrogen.

15. A compound or salt according to any one of Claims 1 to 11 in which

R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

16. A compound or salt according to Claim 15 in which

R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring containing no additional heteroatoms, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_2alkyl, and C_1-C_2alkoxy.

17. A compound or salt according to any one of Claims 2 to 16 wherein

A_1 is aryl, a partially unsaturated heterocyclic group, or heteroaryl group;

substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) –GR_a where

G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

18. A compound or salt according to Claim 17 wherein

A_1 is phenyl, naphthyl, pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, thienylpyrazolyl, benzopyranyl, or 4*H*-chromenyl,

each of which is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, $-CONH_2$, $-SO_2NH_2$, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) $-GR_a$ where

G is chosen from $-(CH_2)_n-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

19. A compound or salt according to Claim 18 in which

A_1 is phenyl, naphthyl, pyridin-2-yl, pyridin-3-yl, pyridin-4-yl, pyrimidin-2-yl, pyrimidinyl-4-yl, pyrimidin-5-yl, thien-2-yl, thien-3-yl, thiazol-4-yl, pyrrol-1-yl, pyrrol-2-yl, pyyrol-3-yl, furan-2-yl, furan-3-yl, pyrazol-1-yl, pyrazol-2-yl pyrazol-4-yl, pyrazol-5-yl, imdiazol-1-yl, imdiazol-2-yl, imdiazol-4-yl, imdiazol-5-yl, thiazol-2-yl, thiazol-3-yl, thiazol-5-yl, *1*,2,3-triazol-4-yl, 1,2,3-thiadiazol-4-yl, 1,2,3-thiadiazol-5-yl, oxazol-2-yl, isoxazol-4-yl, isoxazol-5-yl, isoxazol-3-yl, isoxazol-4-yl, isoxazol-5-yl, benzofuran-2-yl, benzofuran-3-yl, benzopyran-2-yl, benzopyran-3-yl, benzopyran-4-yl, benzo[*d*]oxazol-2-yl benzo[*d*]thiazol-2-yl, benzo[*b*]thiophen-2-yl, 4*H*-chromen-2-yl, benzo[*c*][1,2,5]oxadiazolyl, 2,3-dihydrobenzo[*b*][1,4]dioxin-2-yl, pyrazolo[1,5-a]pyrimidin-6-yl, dihydrobenzo[*b*][1,4]dioxin-3-yl, indol-2-yl, pyrazolo[1,5-*a*]pyrimidin-5-yl, 1H-thieno[2,3-*c*]pyrazol-4-yl, or 1H-thieno[2,3-*c*]pyrazol-5-yl,

each of which is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, $-CONH_2$, $-SO_2NH_2$, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) $-GR_a$ where

G is chosen from $-(CH_2)_n-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

20. A compound or salt according to Claim 19 in which A_1 is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_1-C_6alkoxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_4)alkyl, mono- and di-(C_1-C_4alkyl)amino, and mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl; and

(c) $-GR_a$ where

G is from $-(CH_2)_n-$, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, and

R_a is C_3-C_8cycloalkyl, 5 or 6-membered heterocycloalkyl containing 1 or 2 heteroatoms independently chosen from O, S, and N, 5- or 6-membered heteroaryl containing 1, 2, or 3 heteroatoms independently chosen from O, S, and N, indanyl, and phenyl,

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_2alkyl)amino, and C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

21. A compound or salt according to Claim 20 in which

A_1 is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, C_1-C_6alkyl, C_1-C_6alkoxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_4)alkyl, mono- and di-(C_1-C_4alkyl)amino, and mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl.

22. A compound or salt according to any one of Claim 3 to 16 wherein

A_1 is C_1-C_6 alkyl, C_3-C_7cycloalkyl, or C_2-C_7monocyclic heterocycloalkyl, each of which is substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) –GR_a where

G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

23. A compound or salt according to Claim 22 wherein

A_1 is C_1-C_6 alkyl, C_3-C_7cycloalkyl, pyrrolidinyl, piperidinyl, piperazinyl, or morpholinyl; each of which is substituted with 0 to 3 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

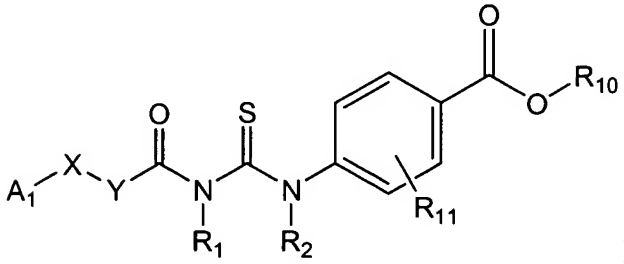
(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) –GR_a where

G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, and phenyl.

24. A compound or salt according to Claim 1 or 2 of Formula 2, wherein :



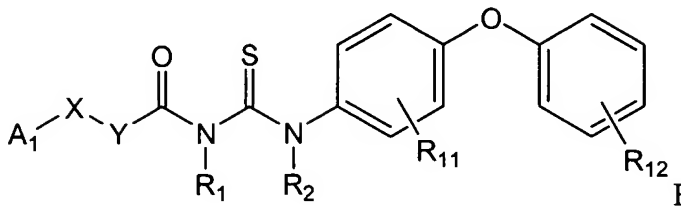
Formula 2

R_1 and R_2 are hydrogen or methyl;

R_{10} is C_1-C_6alkyl; and

R_{11} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

25. A compound or salt according to Claim 1 or 2 of Formula 3, wherein:

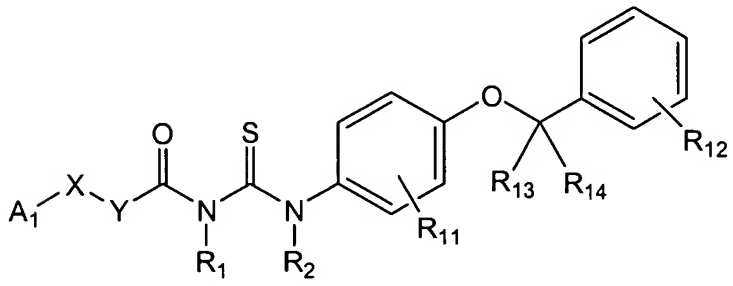


Formula 3

R_1 and R_2 are hydrogen or methyl; and

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

26. A compound or salt according Claim 1 or 2 of Formula 4, wherein:



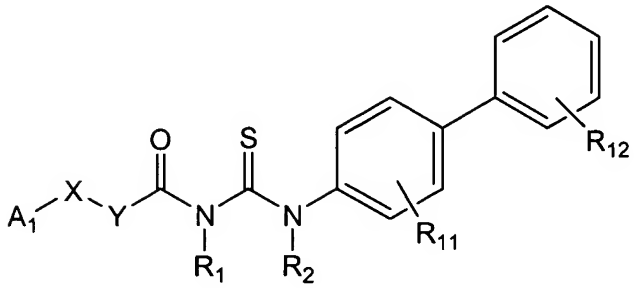
Formula 4

R_1 and R_2 are independently hydrogen or methyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{13} and R_{14} are independently hydrogen or methyl.

27. A compound or salt according to Claim 1 or 2 of Formula 5, wherein

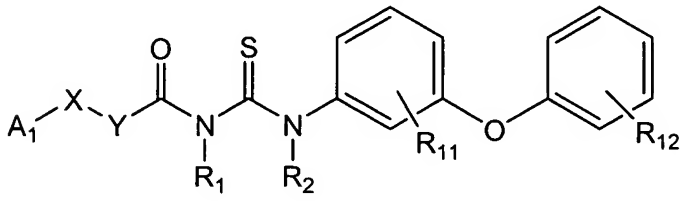


Formula 5

R_1 and R_2 are independently hydrogen or methyl; and

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

28. A compound or salt according to any one of Claim 1 or 2 of Formula 6, wherein

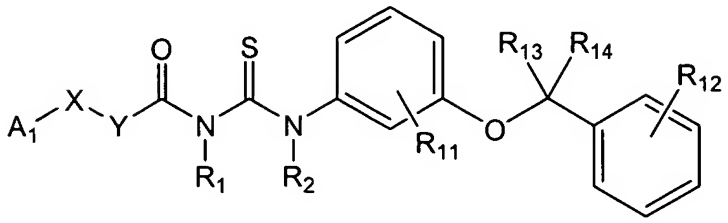


Formula 6

R_1 and R_2 are independently hydrogen or methyl; and

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

29. A compound or salt according to Claim 1 or 2 of Formula 7:



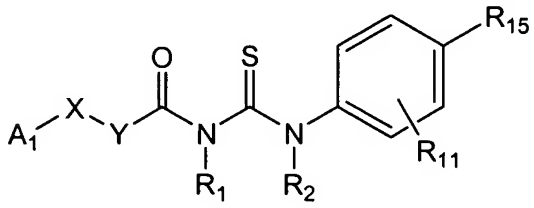
Formula 7

R_1 and R_2 are independently hydrogen or methyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{13} and R_{14} are independently hydrogen or methyl.

30. A compound or salt according to Claim 1 or 2 of Formula 8, wherein



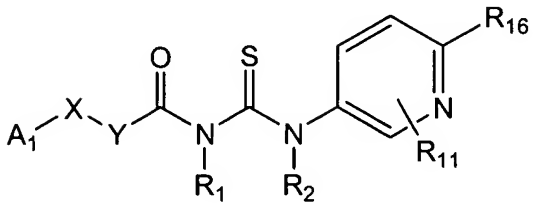
Formula 8

R_1 and R_2 are independently hydrogen or methyl;

R_{11} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{15} represents C_4-C_6alkoxy or C_4-C_6alkyl.

31. A compound salt according to Claim 1 or 2 of Formula 9, wherein



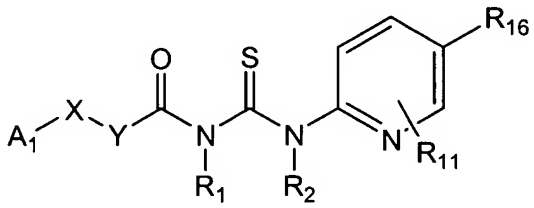
Formula 9

R_1 and R_2 are independently hydrogen or methyl;

R_{11} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{16} is C_2-C_6alkoxy or C_2-C_6alkyl.

32. A compound or salt according to Claim 1 or 2 of Formula 10, wherein



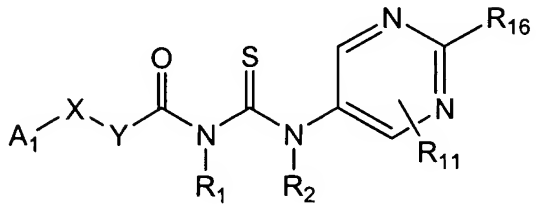
Formula 10

R_1 and R_2 are independently hydrogen or methyl;

R_{11} is represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{16} is C_2-C_6alkoxy or C_2-C_6alkyl.

33. A compound or salt according to Claim 1 or 2 of Formula 11, wherein



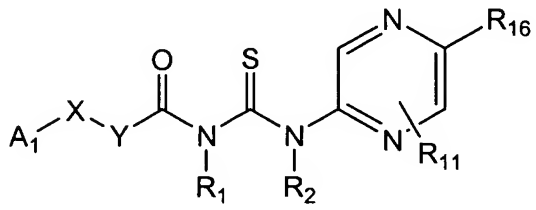
Formula 11

R_1 and R_2 are independently hydrogen or methyl;

R_{11} is represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{16} is C_2-C_6alkoxy or C_1-C_6alkyl.

34. A compound or salt according to Claim 1 or 2 of Formula 12, wherein



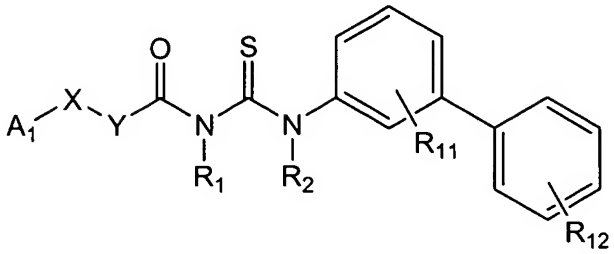
Formula 12

R_1 and R_2 are independently hydrogen or methyl;

R_{11} is represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{16} is C_2-C_6alkoxy or C_2-C_6alkyl.

35. A compound or salt according to Claim 1 or 2 of Formula 13, wherein

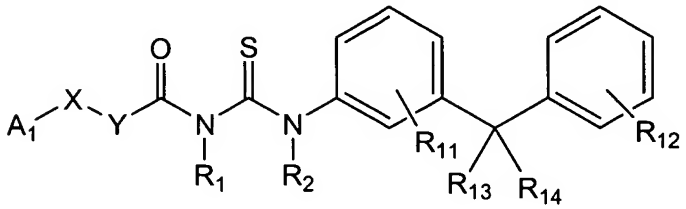


Formula 13

R_1 and R_2 are independently hydrogen or methyl; and

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

36. A compound or salt according to Claim 1 or 2 of Formula 14, wherein



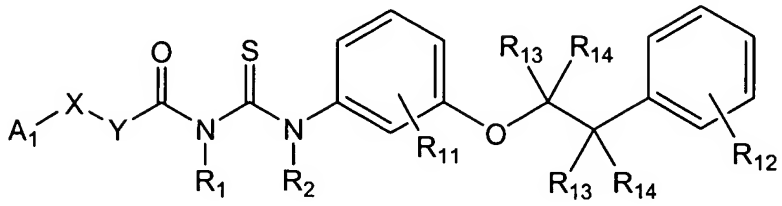
Formula 14

R_1 and R_2 are independently hydrogen or methyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{13} and R_{14} are independently hydrogen or methyl.

37. A compound salt according to Claim 1 or 2 of Formula 15, wherein



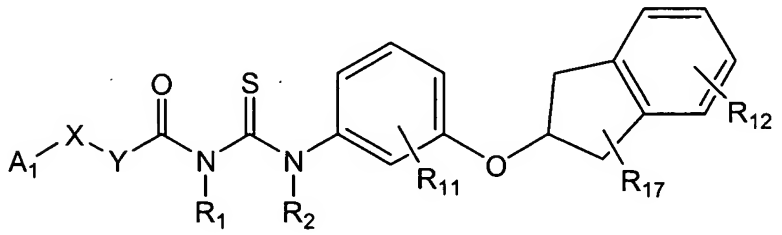
Formula 15

R_1 and R_2 are independently hydrogen or methyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{13} and R_{14} are independently chosen at each occurrence from hydrogen and methyl.

38. A compound salt according to Claim 1 or 2 of Formula 16, wherein



Formula 16

R_1 and R_2 are independently hydrogen or methyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{17} represents 0 to 2 substituents independently chosen from halogen, methyl, and methoxy.

39. A compound or salt according to any one of Claim 24 to 38 in which

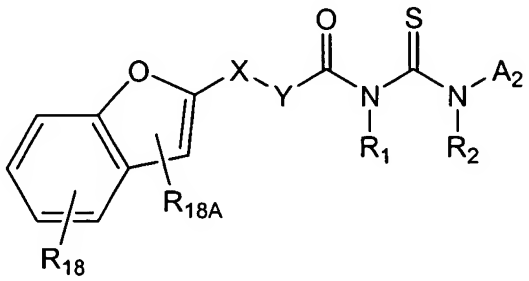
X is NR and Y is $-CH_2-$ or $-CH_2CH_2-$; or

X is O and Y is $-CH_2-$ or $-CH_2CH_2-$; or

X and Y are absent.

40. A compound or salt according to any one of Claims 1, 2, or 24 to 38, wherein A_1 is pyrazinyl, pyridyl, or quinaxolinyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

41. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 17, wherein



Formula 17

wherein

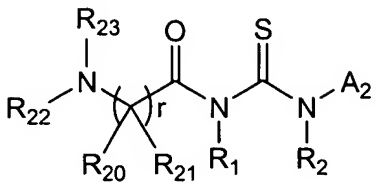
R_{18A} is hydrogen, halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, or C_1-C_2haloalkoxy; and

R_{18} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

42. A compound or salt according to Claim 41 in which

X and Y are absent; and R_1 and R_2 are independently hydrogen or methyl.

43. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 18, wherein



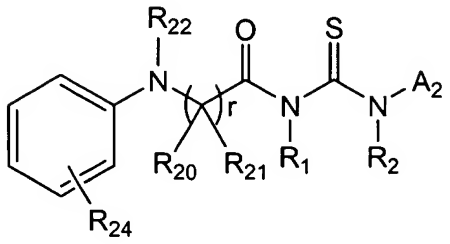
Formula 18

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{22} and R_{23} are independently chosen C_1-C_6 alkyl groups; each of which is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

44. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 19, wherein



Formula 19

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group;

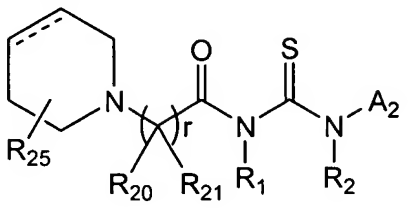
R_{22} is C_1-C_6 alkyl which is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{24} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl.

45. A compound or salt according to Claim 44, wherein

R_{24} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

46. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 20, wherein



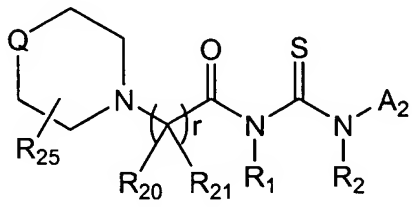
Formula 20

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

47. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 21, wherein



Formula 21

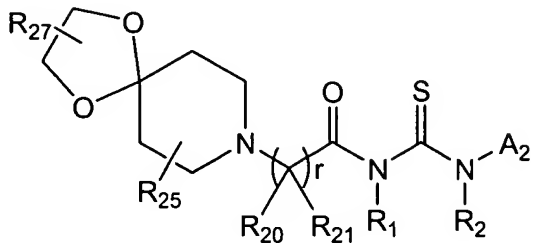
r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

Q is O, S, or NR_{26}; where R_{26} is hydrogen or R_{26} is C_1-C_6alkyl, phenyl, pyridyl, or pyrimidinyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

48. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 22, wherein



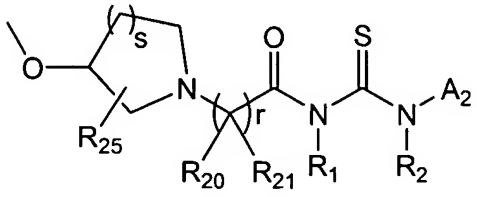
Formula 22

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{25} and R_{27} each represent 0 to 2 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

49. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 23, wherein



Formula 23

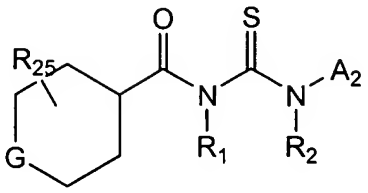
r is 1, 2, or 3;

s is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

50. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 24, wherein



Formula 24

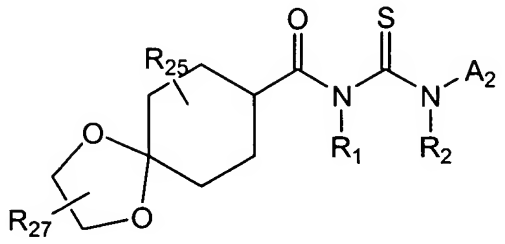
R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

G is O, S, SO_2, or NR_{26}; where R_{26} is hydrogen or R_{26} is C_1-C_6alkyl, phenyl, pyridyl, or pyrimidinyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

51. A compound or salt according to Claim 50, wherein

R_{25} represents a di-(C_1-C_6alkyl)amino substituent and 0 to 2 additional substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

52. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 25, wherein



Formula 25

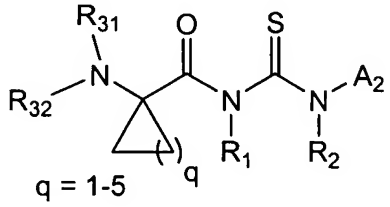
R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{27} represents 0 to 2 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

53. A compound or salt according to Claim 52, wherein

R_{25} represents a di-(C_1-C_6alkyl)amino substituent and 0 to 2 additional substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

54. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 26, wherein

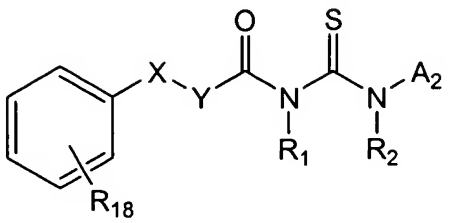


Formula 26

q is an integer from 1 to 5; and

R_{31} and R_{32} are independently chosen from C_1-C_6alkyl and phenyl; each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

55. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 27, wherein



Formula 27

wherein

R_{18} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

56. A compound or salt according to Claim 55 wherein X is oxygen and Y is $-CH_2-$.

57. A compound or salt according to Claim 55 wherein X and Y are absent.

58. A compound or salt according to any one of Claims 1, 2, or 24 to 38, wherein A_1 is 5-membered heteroaryl group selected from furan-2-yl, furan-3-yl, isoxazol-3-yl, isoxazol-4-yl, thiophen-2-yl, thiophen-3-yl, pyrrol-2-yl, pyrrol-3-yl, and pyrazolyl; each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_2-C_4alkenyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

59. A compound or salt according to Claim 58 wherein X is oxygen and Y is $-CH_2-$.

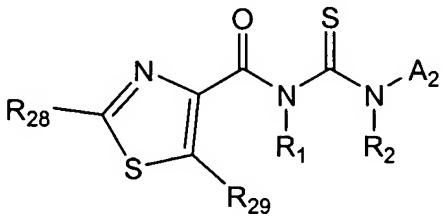
60. A compound or salt according to Claim 58 wherein X and Y are absent.

61. A compound or salt according to any one of Claims 1, 2, or 24 to 38, wherein A_1 is pyridin-2-yl or pyridin-3-yl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

62. A compound or salt according to Claim 61 wherein X is oxygen and Y is $-CH_2-$.

63. A compound or salt according to Claim 61 wherein X and Y are absent.

64. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 28, wherein



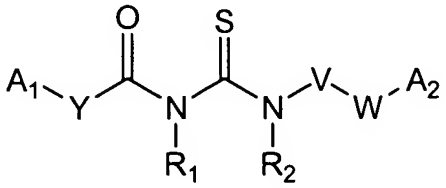
Formula 28

wherein

R_{28} is phenyl or pyridyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{29} is hydrogen, methyl or ethyl.

65. A compound or pharmaceutically acceptable salt of Formula IB, wherein



Formula IB

A_1 is di-(C_1-C_8alkyl)amino, an *N*-(C_1-C_6alkyl)-*N*-phenyl-amino group, an *N*-(C_1-C_6alkyl)-*N*-pyridyl amino group, a 5- to 7-membered monocyclic heterocycloalkyl group covalently bound to a point of attachment in Formula IB via a Nitrogen atom, a 5- to 7-membered monocyclic partially unsaturated heterocyclic group covalently bound to a point of attachment in Formula IB via a Nitrogen atom, a 5- to 7- membered heterocycloalkyl group covalently bound to a point of attachment in Formula IB via a Carbon atom which is adjacent to a Nitrogen atom, or an 8- to 11- membered bicyclic heterocycloalkyl in which the rings are fused or spiro covalently bound to a point of attachment in Formula IB via a Nitrogen atom;

A_2 is C_3-C_8 cycloalkyl, a partially unsaturated or aromatic carbocyclic group, or a saturated, partially unsaturated, or aromatic heterocyclic group;

each of which A_1 and A_2 is substituted with 0 to 5 substituents independently chosen from (a), (b), and (c), where

(a) is independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy,

(b) is independently chosen from C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) is –GR_a where G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl;

W is O, S, NR, or absent, where R is hydrogen or R is C_1-C_6alkyl or aryl(C_0-C_4alkyl), each of which is substituted with 0 to 5 substitutents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, C_1-C_6alkyl, C_1-C_6alkoxy, and mono- and di-(C_1-C_6alkyl)amino;

V is C_1-C_6 alkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent; and when V is absent, W is absent;

Y is C_1-C_6 alkyl substituted with

0 or 1 of C_3-C_7cycloalkyl, a 5- to 7-membered monocyclic heterocycloalkyl, or 8- to 11-membered bicyclic heterocycloalkyl in which the rings are fused or spiro; each of which substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; or

Y is absent;

R_1 and R_2 are independently hydrogen or C_1-C_6alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, or R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

66. A compound or salt according to Claim 65 wherein V and W are absent.

67. A compound or salt according to Claim 65 in which Y is absent.

68. A compound or salt according to Claim 65 in which Y is -CH_2-.

69. A compound or salt according to Claim 65 in which Y is $-CH_2-$ substituted with C_3-C_6cycloalkyl, pyrrolidinyl, or piperidinyl.

70. A compound or salt according to any one of Claims 65 to 69 in which R_1 and R_2 are independently hydrogen or C_1-C_4alkyl.

71. A compound or salt according to Claim 70 in which R_1 and R_2 are independently hydrogen or methyl.

72. A compound or salt according to any one of Claims 65 to 71 wherein

A_2 is C_5-C_7cycloalkyl, phenyl, pyridyl, naphthyl, pyrimidinyl, pyrazinyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently chosen from (a), (b), and (c) where

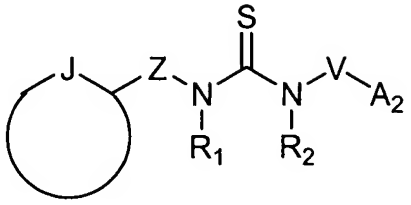
(a) is chosen from halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, $-CONH_2$, $-SO_2NH_2$, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy,

(b) is chosen from C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkanoyl, and C_1-C_8alkoxycarbonyl, and

(c) is $-GR_a$ where G is chosen from $-(CH_2)_n-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and R_a is chosen from C_3-C_8cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

73. A compound of Formula II



Formula II

or a pharmaceutically acceptable salt thereof wherein

A_2 is C_3-C_8 cycloalkyl, a partially unsaturated or aromatic carbocyclic group, a saturated, partially unsaturated, or an aromatic heterocyclic group substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) –GR_a where

G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl;

V is C_1-C_6 alkyl, C_2-C_6alkenyl, or absent; and

Z is carbonyl, thiocarbonyl, or imino;

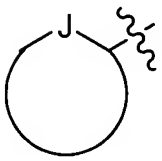
R_1 and R_2 are independently

hydrogen, or

C_1-C_6alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, or

R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy;

the group:



is a group of Formula (i)

that is a saturated, partially unsaturated, or aromatic heterocyclic group where J is O, S, or NR_3 substituted with 0 to 5 substituents independently chosen from: (a), (b), and (c) above; and

R_3 is

(d) hydrogen, C_1-C_2haloalkyl, or C_1-C_2haloalkoxy;

(e) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_4alkoxy(C_1-C_4alkyl), or amino(C_1-C_6)alkyl, or

(f) $-LR_b$ where

L is chosen from $-(CH_2)_r-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_rO(CH_2)_s-$, and $-(CH_2)_rN(CH_2)_s-$, where r and s are independently 0, 1, 2, 3, or 4; and

R_b is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

74. A compound or salt according to Claim 73 wherein Z is carbonyl.

75. A compound or salt according to Claim 73 or Claim 74 wherein V is absent or V is C_1-C_4alkyl.

76. A compound or salt according to any one of Claims 73 to 75 wherein R_1 and R_2 are independently hydrogen or methyl.

77. A compound or salt according to any one of Claims 73 to 76 wherein

A_2 is C_5-C_7cycloalkyl, phenyl, pyridyl, naphthyl, pyrimidinyl, pyrazinyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently chosen from (a), (b), and (c).

78. A compound or salt according to any one of Claims 73 to 77 wherein

A_2 is C_5-C_7cycloalkyl, phenyl, pyridyl, naphthyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, $-CONH_2$, $-SO_2NH_2$, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy,

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkanoyl, and C_1-C_8alkoxycarbonyl, and

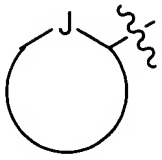
(c) $-GR_a$ where

G is chosen from $-(CH_2)_n-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

79. A compound or salt according to any one of Claims 73 to 78 wherein



is a group of Formula (i)

where Formula (i) is a heteroaryl group that is pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, or thienylpyrazolyl oriented such that the heteroatom J is adjacent to the point of attachment of the group of Formula (i)

the group of Formula (i) is substituted with 0 to 5 substituents independently chosen from: (a), (b), and (c);

J is S, O, or NR_3; and

R_3 is

(d) hydrogen, C_1-C_2haloalkyl, or C_1-C_2haloalkoxy;

(e) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_4alkoxy(C_1-C_4alkyl), or amino(C_1-C_6)alkyl, or

(f) $-LR_b$ where

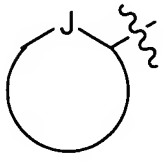
L is chosen from $-(CH_2)_r-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_rO(CH_2)_s-$, and $-(CH_2)_rN(CH_2)_s-$, where r and s are independently 0, 1, 2, 3, or 4; and

R_b is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

80. A compound or salt of Claim 79 wherein

wherein



is a group of Formula (i)

where Formula (i) is a heteroaryl group that is pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, or thienylpyrazolyl oriented such that the heteroatom J is adjacent to the point of attachment of the group of Formula (i) in Formula II;

the group of Formula (i) is substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, $-CONH_2$, $-SO_2NH_2$, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkanoyl, and C_1-C_8alkoxycarbonyl,

(c) $-GR_a$ where

G is chosen from $-(CH_2)_n-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl;

J is S, O, or NR_3, and

R_3 is

(d) hydrogen,

(e) C_1-C_6alkyl, or

(f) $-LR_b$ where

L is chosen from $-(CH_2)_r-$, $-(CH_2)_rO(CH_2)_s-$, and $-(CH_2)_rN(CH_2)_s-$, where r and s are independently 0, 1, 2, 3, or 4; and

R_b is chosen from C_3-C_8cycloalkyl, piperidinyl, piperazinyl, morpholinyl, indanyl, tetrahydronapthyl, phenyl, and pyridyl;

each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

81. A compound or pharmaceutically acceptable salt thereof, in which the compound is selected from

1-(Furan-2-carbonyl)-3-(4-benzo[*d*]thiazol-2-yl-phenyl)-thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-[5-(benzo[*d*]oxazol-2-yl)-2-methyl]phenylthiourea;

1-(3-(Benzo[*d*]thiazol-2-yl)phenyl)-3-(2-phenoxyacetyl) thiourea;

1-(4-(Benzo[*d*]oxazol-2-yl)phenyl)-3-propionylthiourea;

1-(Pyridin-3-carbonyl)-3-(4-benzo[*d*]thiazol-2-yl-phenyl)-thiourea;

1-[3-(2-chlorophenyl-5-methyl-isoxazol-4-yl)-carbonyl]-3-(4-isopropylphenyl)thiourea;

Butyl4-(3-(2-phenoxyacetyl) thioureido)benzoate;

Butyl 4-(3-acetylthioureido)benzoate;

Butyl 4-(3-(2-(3-chlorophenoxy) acetyl) thioureido)benzoate;

Butyl 4-(3-(3-phenoxypropanoyl) thioureido)benzoate;

Butyl 4-(3-(2-(naphthalen-3-yloxy)acetyl)thioureido)benzoate;

Butyl 4-(3-(benzofuran-2-yl-carbonyl)thioureido)benzoate;

Ethyl 2-(4-(3-(2-phenoxyacetyl)thioureido)phenyl)acetate;

Ethyl 4-(3-(2-phenoxyacetyl)thioureido)benzoate;

Butyl 4-(3-(2-methoxyacetyl) thioureido)benzoate;

Butyl 4-(3-(2-(2,4-dichlorophenoxy) acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(4-tert-butylphenoxy) acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(4-(benzyloxy) phenoxy)acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(2-methoxyphenoxy) acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(o-tolyloxy)acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(2,4,6-trichlorophenoxy)acetyl)thioureido)benzoate;

Butyl 4-(3-(3,4-dichlorophenyl) carbonyl)thioureido)benzoate;

1-(3,4-dichlorophenyl-carbonyl)-3-(3-triflurомethylphenyl)thiourea;

1-(3,4-Dichlorophenyl-carbonyl)-3-(3-benzoxy-phenyl)thiourea;

1-(3,4-Dichlorophenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3,4-Diflurophenyl-carbonyl)-3-(3-(5-methylfuran-2-yl)-phenyl)thiourea;

Butyl 4-(3-(naphth-2-yl) carbonyl)thioureido)benzoate;

Butyl 4-(3-(4-cyanophenyl) carbonyl)thioureido)benzoate;

Butyl 4-(3-(methylacetate) carbonyl)thioureido)benzoate;

1-((Benzofuran-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;

Butyl 4-(3-(2-(3,4-dichlorophenoxy)acetyl)thioureido)benzoate;

1-(4-Butylphenyl)-3-(2-phenoxyacetyl)thiourea;

(Amino-(3-(benzyloxy)phenyl) methanethiocarbamoyl)methyl acetate;

1-(3-(Methylthio)propanoyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(2-(2-Chlorophenoxy)acetyl)-3-(3-(benzyloxy)phenyl)thiourea;

Butyl 4-(3-(naphth-1-yl) carbonyl)thioureido)benzoate;

(S)-1-(Amino-*N*-p-(butylacetate) methanethiocarbamoyl)ethyl acetate;

Butyl 4-(3-(2-(2-methoxyethoxy)acetyl)thioureido)benzoate;

(Amino*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)(phenyl)methyl acetate;

Ethyl 3-(amino-*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)propanoate;

1-Butyryl-3-(4-cyclohexylphenyl)thiourea;

(S)-1-(Amino-*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)ethyl acetate;

1-(3-(Benzyloxy)phenyl)-3-(2-hydroxyacetyl)thiourea;

Butyl 4-(3-(2-(2,6-dichlorophenoxy) acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(3-methoxyphenoxy) acetyl) thioureido)benzoate;

1-[(1-methyimidazol-2-yl) –carbonyl]-3-(3-benzoxy-phenyl)thiourea;

tert-Butyl 2-(aminoN-(3-(benzyloxy) phenyl)methanethiocarbamoyl) pyrrolidone-1-carboxylate;

Butyl 4-(3-(pyrrolidin-1-yl) carbonyl)thioureido)benzoate;

Butyl 4-(3-(1-methyl-benzofuran-2-yl) carbonyl)thioureido)benzoate;

1-(4-Hexylphenyl)-3-(2-phenoxyacetyl)thiourea;

1-(4-(Pentyloxy)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyloxy)-phenyl)thiourea;

1-(4-Pentylphenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyl)-phenyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyloxy)-phenyl)thiourea;

1-(4-Butoxyphenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-phenyl)-phenyl)thiourea;

1-(2-Phenoxyacetyl)-3-(3-phenyl)-phenylthiourea;

Isopropyl 4-(3-(benzofuran-2-yl) carbonyl)thioureido)benzoate;

1-(2-phenoxyacetyl)-3-94-fluoro-phenyl)-phenylthiourea;

1-(3-benzylphenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-benzyl)-phenyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-benzyl)-phenyl)thiourea;

1-(4-(p-Tolyloxy)phenyl)-3-(2-phenoxyacetyl)thiourea;

Isobutyl 4-(3-(benzofuran-2-yl) carbonyl)thioureido)benzoate;

Isobutyl 4-(3-(2-phenoxyacetyl)thioureido)benzoate;

1-(2-(phenylmethanone)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-(Phenylcarbamoyl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-(2-Methylpyrimidin-4-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(4-(4-Chlorophenoxy)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-Chlorophenoxy)-phenyl)thiourea;

1-(4-(3,4-Dihydroisoquinolin-2(1H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-Fluoro-4-(octahydroquinolin-1(2H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-Fluoro-4-(octahydroquinolin-1(2H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-Fluoro-4-(piperidin-1-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-fluoro-4-(piperidin-1-yl)phenyl)-phenyl)thiourea;

1-(3-(3-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-(3-methoxybenzyloxy)) phenyl)-phenyl)thiourea;

1-(3-(2-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-(2-Methoxybenzyloxy) phenyl)-thiourea;

1-(3-(4-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-(cyclohexylmethoxy) phenyl)-thiourea;

1-(3-(Cyclohexylmethoxy) phenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-(5,6-dihydropyridin-1(2H)-yl))phenyl)-thiourea;

1-((5-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-benzyloxy-phenyl)-thiourea;

Butyl 4-(3-(5-chloro-benzofuran-2-yl) carbonyl)thioureido)benzoate;

1-(7-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(2,3,4-tetrahydronapthalen-2-yl-carbonyl)-3-(3-(3-methoxybenzyloxy))phenyl)-phenyl)thiourea;

1-(2-(4-(Trifluoromethoxy) phenoxy) acetyl)-3-(3-(benzyloxy)phenyl) thiourea;

1-(3-(Benzyloxy)phenyl)-3-(2-(pyridin-3-yloxy)acetyl) thiourea;

1-(4-oxo-4-*H*-chromen-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(3-(Benzyloxy)phenyl)-3-(2-(pyridin-2-yloxy)acetyl) thiourea;

1-(Pyridin-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(3-Chloro-benzo[*b*]thiophen-2-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(4-Trifluoromethoxy-phenyl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(5-Methylisoxazol-3-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(2-Methyl-5-phenyl-furan-3-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(4-Trifluoromethyl-phenyl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(3-Chloro-benzo[b]thiophen-2-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(3,5-Dimethylisoxazol-4-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(5-Methylisoxazol-3-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(4-Trifluoromethylphenyl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

(S)-1-(Amino-*N*-(3-(benzyloxy) phenyl)methanethiocarbamoyl)ethyl acetate;

(S)-1-(Amino-*N*-(3-phenoxyphenyl)methane thiocarbamoyl)ethyl acetate;

(S)-1-(Amino-*N*-(3-benzyl-phenyl)methanethiocarbamoyl)ethyl acetate;

Ethyl 1-(2-fluoro-4-(3-(benzofuran-2-yl-carbonyl)thioureido)phenyl)-4-phenylpiperidine-4-carboxylate;

1-(4-Trifluoromethylphenyl-carbonyl)-3-(2-phenylbenzo[d][1,3]dioxol-6-yl) phenyl)thiourea;

1-(3-Chloro-methylbenzo[b]thiophen-2-yl-carbonyl)-3-(2-phenylbenzo[d][1,3]dioxol-6-yl) phenyl)thiourea;

1-(4-Trifluoromethoxyphenyl-carbonyl)-3(2-phenylbenzo[d][1,3]dioxol-6-yl) thiourea;

1-(5-Methylisoxazol-3-yl-carbonyl)-3-(2-phenylbenzo[d][1,3]dioxol-6-yl) phenyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(3-chloro-methylbenzo[b]thiophen-2-yl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethoxy-phenyl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)thiourea;

1-(3-((S)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)-thiourea;

1-(3-((S)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)-thiourea;

1-(3-((S)-1-Phenylethoxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(3-chloro-methylbenzo[b]thiophen-2-yl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(3,5-dimethylisoxazol-4-yl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)-thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(3,5-dimethylisoxazol-4-yl-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)thiourea;

1-(2-Phenylbenzo[*d*][1,3]dioxol-6-yl)-3-(benzofuran-2-yl-carbonyl)thiourea;

1-(3-((S)-1-Phenylethoxy)phenyl)-3-(benzofuran-2-yl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(benzofuran-2-yl-carbonyl)-thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(benzofuran-2-yl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(benzofuran-2-yl-carbonyl)thiourea;

1-(2,4-dimethylthiazol-5-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(1-methyl-pyrrol-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(3-(Phenethyloxy)phenyl)-3-((2,7-dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl)-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(1-ethyl-3-methyl-1*H*-pyrazol-5-yl-carbonyl)-thiourea;

1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-(2,4-dimethylthiazol-5-yl-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-((2,7-dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl)-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-(1-ethyl-3-methyl-1*H*-pyrazol-5-yl-carbonyl)thiourea;

1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(Benzo[*d*]thiazol-2-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(Benzo[*d*]thiazol-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-fluoro-5-pentoxy-phenyl)thiourea;

1-(2-Methyl-pyridin-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(1-phenyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(1-phenyl-1*H*-pyrazol-5-yl)-3-((3-phenyloxy)phenyl)-phenyl)thiourea;

1-(2-Phenylbenzo[*d*][1,3]dioxol-6-yl)-3-(1-phenyl-1*H*-pyrazol-5-yl-carbonyl)thiourea;

1-(1-phenyl-1*H*-pyrazol-5-yl)-3-(4-pentoxy-phenyl)-phenyl)thiourea;

1-(1-phenyl-1*H*-pyrazol-5-yl)-3-((3-phenyloxy-phenyl)-phenyl)thiourea;

1-(Methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-((4-pentoxy-phenyl)-phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(4-(pentyl) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(4-(pentoxy) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-((3-(trifluoromethyl)phenyl)furan-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-((3-(trifluoromethyl)phenyl)furan-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(5-Nitro-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(5-Nitro-benzofuran-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(4-pentyl phenyl)thiourea;

1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(4-pentoxy phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-(hex-1-ynyl)phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethynyl)phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethyl)phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(pyridin-2-yl) phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(pyridin-2-yl-oxy)-phenyl) thiourea;

1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(4-(hex-1-ynyl)phenyl) thiourea;

1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethynyl)phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-((3-(trifluoromethyl)benzyloxy)-phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-((1-methylpiperidin-4-yl)methoxy)-3-fluorophenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(trifluoromethyl)-4-(piperidin-1-yl)phenyl)thiourea;

1-(1,3-dimethyl-1*H*-thieno[2,3-*c*]pyrazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(4-(pentyl) phenyl)thiourea;

1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(4-(pentoxy) phenyl)thiourea;

1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;

1-(3-Methylisoxazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(3-Methylisoxazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(3-Methylisoxazol-4-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;

1-(3-Methylisoxazol-4-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(5-Methyl-2-phenyl-2*H*-1,2,3-triazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(5-Methyl-2-phenyl-2*H*-1,2,3-triazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(5-Chloro-3-methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(5-Chloro-3-methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;

1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(3-(phenyloxy)-phenyl)thiourea;

1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;

1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(4-Methoxy-benzofuran-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;

1-(4-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-fluoro-4-pentoxy-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3,5-dibromo-4-(pent-4-enyloxy)phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-((pyridin-3-yl)methyl) phenyl)thiourea;

1-(5-Iodo-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(5-Phenyl-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(5-(2-Pyridyl)benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(3-Propoxy-pyridin-2-yl-carbonyl)-3-(4-(pentyl)phenyl)thiourea;

1-(2,5-Dichlorothiophen-3-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;

1-(5-(Methylthio)-thiophen-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;

1-(5,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;

1-(2,5-Dichlorothiophen-3-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;

1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;

1-(5-(Methylthio)-thiophen-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;

1-(7-Fluoro-benzofuran-2-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(7-Fluoro-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(2-(1,3-Dioxoisoindolin-2-yl)acetyl)-3-(3-phenoxyphenyl)thiourea;

1-(2-(1,3-Dioxoisoindolin-2-yl)acetyl)-3-(3-benzyloxyphenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)methyl-phenyl)thiourea ;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(phenylamino)methyl-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-(*N*-benzyl-N-methylamino)-3-fluorophenyl) thiourea;

Phenyl 3-(3-((benzofuran-2-yl)-carbonyl)thioureido)benzoate;

1-(4-Cyanophenyl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(4-Cyanophenyl-carbonyl)-3-(4-(pentyloxy)-phenyl)thiourea;

1-(4-Cyanophenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(Quinoxalin-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(Quinoxalin-2-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(Quinoxalin-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(4-Trifluoromethylphenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(4-Trifluoromethylphenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(5-Cyano-benzofuran-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(5-Cyano-benzofuran-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

Butyl 4-(3-(4-Methyl-1,2,3-thiadiazol-5-yl-) carbonyl)thioureido) benzoate;

1-(Pyrazin-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(Pyrazin-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(Pyrazin-2-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Fluoro-4-(pentyloxy)phenyl)-3-(2-(1,3-dioxoisoindolin-2-yl)acetyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-trifluoromethyl-4-pentoxy-phenyl)thiourea;

1-(1-Benzyl-1*H*-tetrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-(N-methyl-N-pentylamino)-3-fluorophenyl)thiourea;

1-(1-Benzyl-1*H*-tetrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(6-Trifluoromethyl-pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(6-Trifluoromethyl-pyrid-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Trifluoromethoxy-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Trifluoromethoxy-phenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(2-Chloro-5-trifluoromethoxy-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Difluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Difluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(5-(Trifluoromethyl)-2-phenyloxazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(5-(Trifluoromethyl)-2-phenyloxazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(5-(2-Chloro-5-trifluoromethylphenyl)-furan-2-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(3-Trifluoromethyl-4-methoxy-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-4-chloro-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Trifluoromethyl-4-chloro-phenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea;

1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-((5-Acetamidobenzofuran-2-yl)carbonyl)-3-(3-phenoxyphenyl)thiourea;

1-Acetyl-3-(3-phenoxyphenyl)thiourea;

1-Acetyl-3-(4-(pentyloxy)phenyl)thiourea;

1-Acetyl-3-(4-pentylphenyl)thiourea;

1-(Dimethylamino-acetyl)-3-(3-phenoxyphenyl)thiourea;

1-(Dimethylamino-acetyl)-3-(3-benzyloxyphenyl)thiourea;

1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-((2,3,4,5,6-penta-fluorophenoxy)-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(9-methyl-9*H*-fluoren-7-yl)thiourea;

Pentyl 2-phenyl- 4-(3-(benzofuran-2-yl)thioureido)benzoate;

1-(3-Pyrid-3-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea;

1-(3-Pyrid-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(4-Phenylbutanoyl)-3-(3-phenoxyphenyl)thiourea;

1-(4-Phenylbutanoyl)-3-(3-benzyloxyphenyl)thiourea;

1-(2-Morpholinoacetyl)-3-(3-phenoxyphenyl)thiourea;

1-(2-Morpholinoacetyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(2-Morpholinoacetyl)-3-(4-(pentyl)phenyl)thiourea;

1-(4-(Pentyloxy)phenyl)-3-(2-(piperidin-1-yl)acetyl)thiourea;

1-(*N*-Methyl-*N*-phenylamino-acetyl)-3-(3-benzyloxyphenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(6-pentoxy-pyrid-3-yl)thiourea;

1-(3-Pyrid-3-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea hydrochloride;

1-(3-Pyrid-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea hydrochloride;

1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea hydrochloride;

1-(Benzofuran-2-yl-carbonyl)-3-(trifluoromethylthio-phenyl)thiourea;

1-(3-(Piperidin-1-yl)propanoyl)-3-(4-pentylphenyl)thiourea;

1-(3-(Piperidin-1-yl)propanoyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(3-(Piperidin-1-yl)propanoyl)-3-(3-phenoxyphenyl)thiourea;

1-(3-Morpholinopropanoyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(1-Methylpiperidin-3-yl-carbonyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(1-Methylpiperidin-3-yl-carbonyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(2-(2-methylpiperidin-1-yl)acetyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(2-Oxo-4-phenyl-pyrrolidin-1-ylcarbonyl)-3-(3-benzyloxy-phenyl)thiourea; and

1-(5-Trifluoromethoxy-benzofuran-2-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea.

82. A pharmaceutical composition comprising a compound or salt according to any one of Claims 1 to 81 together with a pharmaceutically acceptable carrier, diluent, or excipient.

83. A pharmaceutical composition according to Claim 82, wherein the composition is formulated as an injectable fluid, an aerosol, a cream, a gel, a pill, a capsule, a syrup, a tablet, ophthalmic solution, or a transdermal patch.

84. A package comprising a pharmaceutical composition of Claim 82 in a container and further comprising instructions for using the composition to treat a patient suffering from Hepatitis C infection.

85. A compound or salt according to Claim 1 that exhibits an EC_{50} of less than 10 micromolar in a replicon assay of HCV replication.

86. A compound or salt according to Claim 1 that exhibits an EC_{50} of less than 1 micromolar in a replicon assay of HCV replication.

87. A method for treating Hepatitis C infection comprising administering to a patient in need of such treatment a therapeutically effective amount of a compound or salt according to Claim 1.

88. The method of Claim 87 wherein the patient is a human patient.

89. The method of Claim 88 wherein the therapeutically effective amount is an amount sufficient to significantly decrease the number of HCV antibodies in the patient's blood or serum.

90. A method of inhibiting HCV replication *in vivo* comprising administering to a patient infected with HCV a concentration of a compound or salt according to Claim 1 sufficient to inhibit HCV replicon replication *in vitro*.